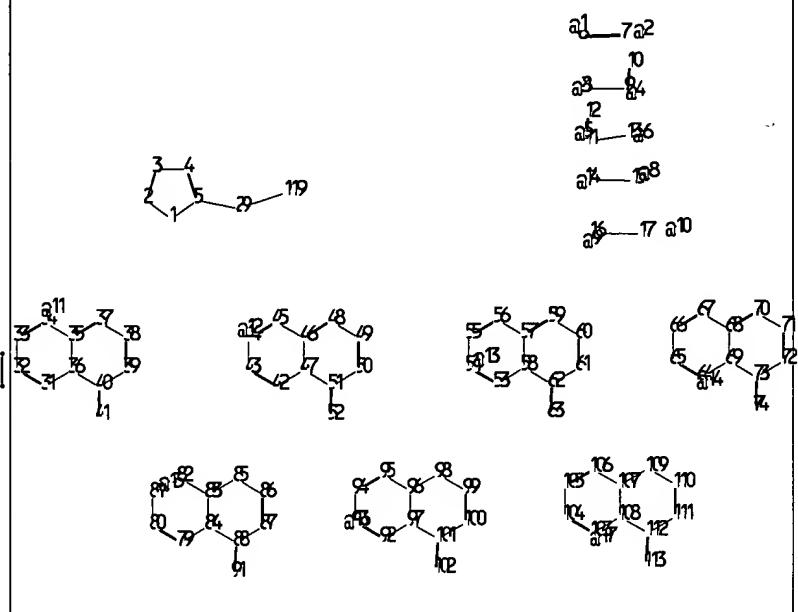
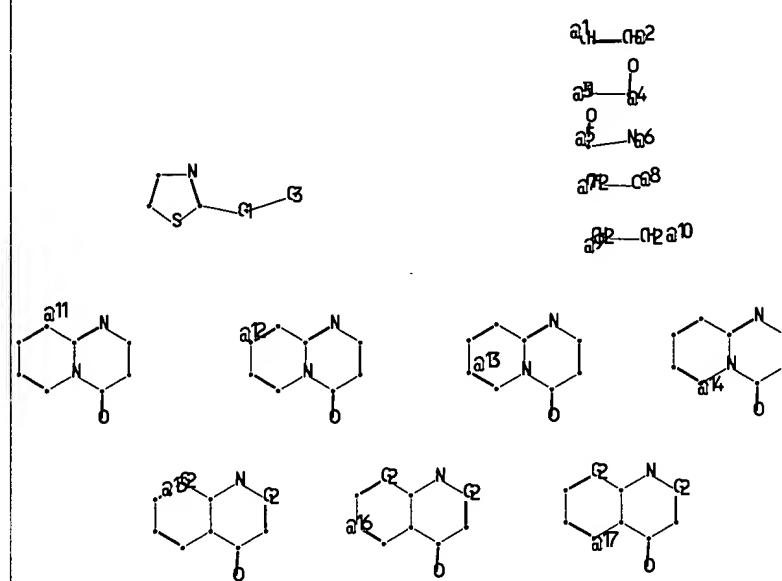


L Number	Hits	Search Text	DB	Time stamp
1	3144	((514/233.2) or (514/248) or (514/258) or (514/300) or (514/312)).CCLS.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2002/02/11 11:26
2	3272	((544/116) or (544/235) or (544/236) or (544/282) or (546/122) or (546/153) or (546/155) or (546/156)).CCLS.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2002/02/11 11:27
3	5449	((((514/233.2) or (514/248) or (514/258) or (514/300) or (514/312)).CCLS.) or ((544/116) or (544/235) or (544/236) or (544/282) or (546/122) or (546/153) or (546/155) or (546/156)).CCLS.)	USPAT; US-PGPUB; EPO; JPO; DERWENT	2002/02/11 11:27
5	38047	thiazolyl or thiazole	USPAT; US-PGPUB; EPO; JPO; DERWENT	2002/02/11 11:28
6	1229	((((514/233.2) or (514/248) or (514/258) or (514/300) or (514/312)).CCLS.) or ((544/116) or (544/235) or (544/236) or (544/282) or (546/122) or (546/153) or (546/155) or (546/156)).CCLS.)) and (thiazolyl or thiazole)	USPAT; US-PGPUB; EPO; JPO; DERWENT	2002/02/11 11:28



chain nodes :

6 7 8 9 10 11 12 13 14 15 16 17 29 41 52 63 74 91 102 113 119

ring nodes :

106 107
chain bonds :

In bonds : 5-29 6-7 8-9 9-10 11-12 11-13 14-15 16-17 29-119 40-41 51-52 62-63 73-74 88-91
 101-102 112-113

ring bonds :

exact/norm bonds :

107-108 107-109 1 isolated ring systems :

containing 1 : 31 : 42 : 53 : 64 : 79 : 82 : 103 :

G2:C,N

G3:[*11], [*12], [*13], [*14], [*15], [*16], [*17]

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:CLASS 7:CLASS 8:CLASS 9:CLASS 10:CLASS
11:CLASS 12:CLASS 13:CLASS 14:CLASS 15:CLASS 16:CLASS 17:CLASS 29:CLASS 31:Atom
32:Atom 33:Atom 34:Atom 35:Atom 36:Atom 37:Atom 38:Atom 39:Atom 40:Atom 41:CLASS
42:Atom 43:Atom 44:Atom 45:Atom 46:Atom 47:Atom 48:Atom 49:Atom 50:Atom 51:Atom
52:CLASS 53:Atom 54:Atom 55:Atom 56:Atom 57:Atom 58:Atom 59:Atom 60:Atom 61:Atom
62:Atom 63:CLASS 64:Atom 65:Atom 66:Atom 67:Atom 68:Atom 69:Atom 70:Atom 71:Atom
72:Atom 73:Atom 74:CLASS 79:Atom 80:Atom 81:Atom 82:Atom 83:Atom 84:Atom 85:Atom
86:Atom 87:Atom 88:Atom 91:CLASS 92:Atom 93:Atom 94:Atom 95:Atom 96:Atom 97:Atom
98:Atom 99:Atom 100:Atom 101:Atom 102:CLASS 103:Atom 104:Atom 105:Atom 106:Atom
107:Atom 108:Atom 109:Atom 110:Atom 111:Atom 112:Atom 113:CLASS 119:CLASS

09/842,234

=>
Uploading 09842234.str

L1 STRUCTURE UPLOADED

=> d 11
L1 HAS NO ANSWERS
L1 STR
*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

Structure attributes must be viewed using STN Express query preparation.

=> s 11 sss sam
SAMPLE SEARCH INITIATED 09:43:14 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 44 TO ITERATE

100.0% PROCESSED 44 ITERATIONS 1 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 483 TO 1277
PROJECTED ANSWERS: 1 TO 80

L2 1 SEA SSS SAM L1

=> s 11 sss ful
FULL SEARCH INITIATED 09:43:38 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 1000 TO ITERATE

100.0% PROCESSED 1000 ITERATIONS 22 ANSWERS
SEARCH TIME: 00.00.05

L3 22 SEA SSS FUL L1

=> s 13
L4 2 L3

=> d 14 1-2 bib,ab,hitstr

L4 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2002 ACS
 AN 2001:319871 CAPLUS
 DN 134:336205
 TI Drug discharge pump inhibitors
 IN Leger, Roger; Watkins, William John; Zhang, Jason Zhijia; Renau, Thomas Eric; Lee, Ving Jack; Ohta, Toshiharu; Nakayama, Kiyoshi; Ishida, Yohhei; Ohtsuka, Masami; Kawato, Haruko
 PA Microcide Pharmaceuticals, Inc., USA; Daiichi Pharmaceutical Co., Ltd.
 SO PCT Int. Appl., 237 pp.
 CODEN: PIXXD2

DT Patent

LA Japanese

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2001030757	A1	20010503	WO 2000-JP7565	20001027
	W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NC, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
	RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				

PRAI US 1999-428466 A 19991028
 JP 2000-326713 A 20001026

AB Drugs for preventing and/or treating microbial infectious diseases which contain, as the active ingredient, compds. represented by the formula R1R2J1W1A1(G1)m[CH(R3)]p(G2)nG3Q1, physiol. acceptable salts thereof or hydrates of the same and have an effect of making a microorganism having acquired tolerance to a drug non-tolerant. In said formula R1 and R2 independently represent each hydrogen, halogeno, carboxy, etc.; J1 represents 5- or 6-membered heteroaryl; W1 represents -CH=CH-, -CH CH-, -CH2CH2-, etc.; A1 represents phenylene, pyridinedyl, furandyl, etc.; G1 represents oxygen, carbonyl, ethynyl, etc.; p is an integer of from 0 to 3; G2 represents phenylene, furandyl, tetrahydrofurandyl, etc.; G3 represents -CH2- or a single bond; m and n represent each an integer of 0 or 1; and Q1 represents an acidic group.

IT 337904-40-2P 337904-41-3P 337904-43-5P
 337904-45-7P 337904-46-8P 337904-47-9P
 337904-48-0P 337904-49-1P 337904-50-4P
 337904-51-5P 337904-52-6P 337904-53-7P
 337904-56-0P 337904-57-1P 337904-58-2P
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 337904-62-8P 337904-63-9P 337904-64-0P

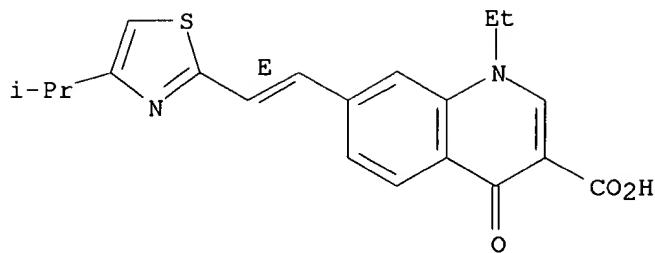
RL: BAC (Biological activity or effector, except adverse); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(drug discharge pump inhibitors as antimicrobials)

RN 337904-40-2 CAPLUS

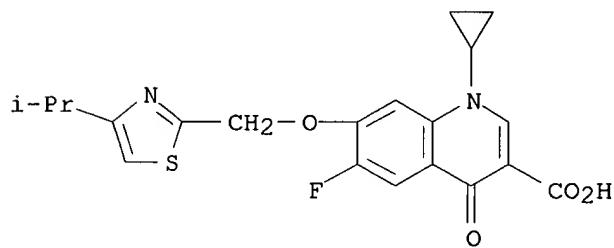
CN 3-Quinolinecarboxylic acid, 1-ethyl-1,4-dihydro-7-[(1E)-2-[4-(1-methylethyl)-2-thiazolyl]ethenyl]-4-oxo- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 337904-41-3 CAPLUS

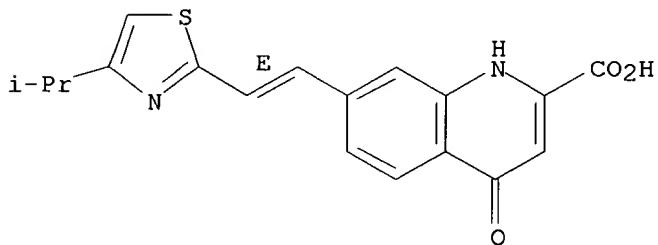
CN 3-Quinolinecarboxylic acid, 1-cyclopropyl-6-fluoro-1,4-dihydro-7-[(4-(1-methylethyl)-2-thiazolyl)methoxy]-4-oxo- (9CI) (CA INDEX NAME)



RN 337904-43-5 CAPLUS

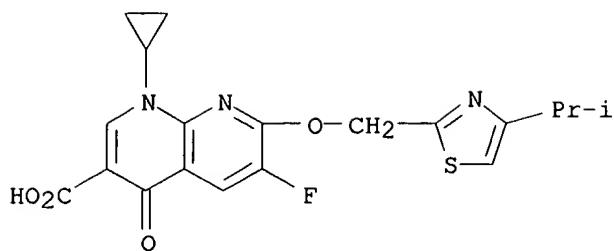
CN 2-Quinolinecarboxylic acid, 1,4-dihydro-7-[(1E)-2-[(4-(1-methylethyl)-2-thiazolyl)ethenyl]-4-oxo- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 337904-45-7 CAPLUS

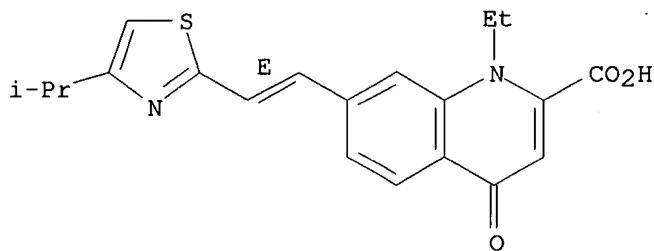
CN 1,8-Naphthyridine-3-carboxylic acid, 1-cyclopropyl-6-fluoro-1,4-dihydro-7-[(4-(1-methylethyl)-2-thiazolyl)methoxy]-4-oxo- (9CI) (CA INDEX NAME)



RN 337904-46-8 CAPLUS

CN 2-Quinolinecarboxylic acid, 1-ethyl-1,4-dihydro-7-[(1E)-2-[4-(1-methylethyl)-2-thiazolyl]ethenyl]-4-oxo- (9CI) (CA INDEX NAME)

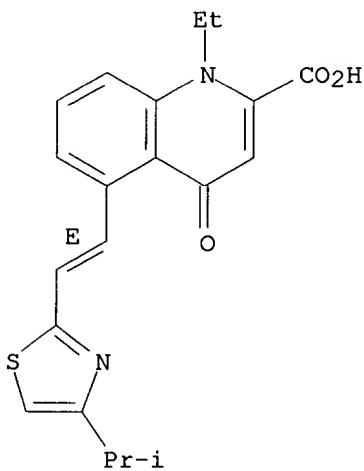
Double bond geometry as shown.



RN 337904-47-9 CAPLUS

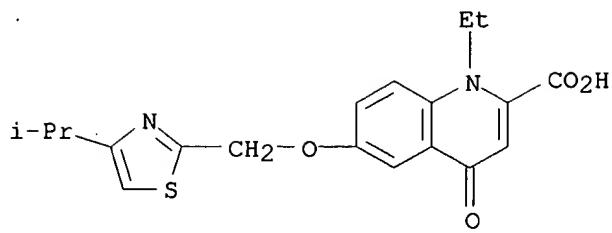
CN 2-Quinolinecarboxylic acid, 1-ethyl-1,4-dihydro-5-[(1E)-2-[4-(1-methylethyl)-2-thiazolyl]ethenyl]-4-oxo- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



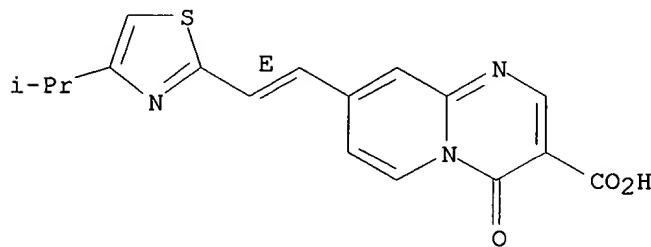
RN 337904-48-0 CAPLUS

CN 2-Quinolinecarboxylic acid, 1-ethyl-1,4-dihydro-6-[(4-(1-methylethyl)-2-thiazolyl)methoxy]-4-oxo- (9CI) (CA INDEX NAME)

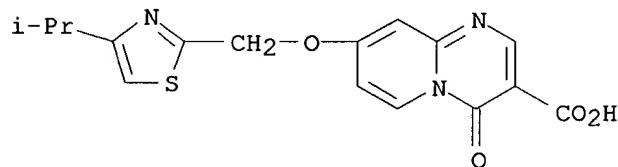


RN 337904-49-1 CAPLUS
 CN 4H-Pyrido[1,2-a]pyrimidine-3-carboxylic acid, 8-[(1E)-2-[4-(1-methylethyl)-2-thiazolyl]ethenyl]-4-oxo- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

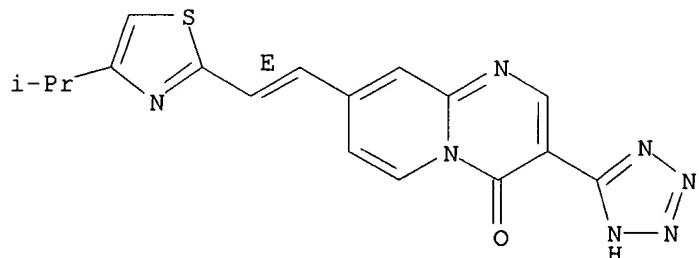


RN 337904-50-4 CAPLUS
 CN 4H-Pyrido[1,2-a]pyrimidine-3-carboxylic acid, 8-[(4-(1-methylethyl)-2-thiazolyl)methoxy]-4-oxo- (9CI) (CA INDEX NAME)



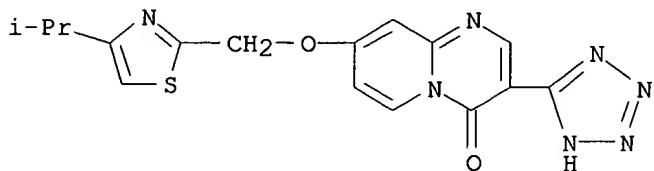
RN 337904-51-5 CAPLUS
 CN 4H-Pyrido[1,2-a]pyrimidin-4-one, 8-[(1E)-2-[4-(1-methylethyl)-2-thiazolyl]ethenyl]-3-(1H-tetrazol-5-yl)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



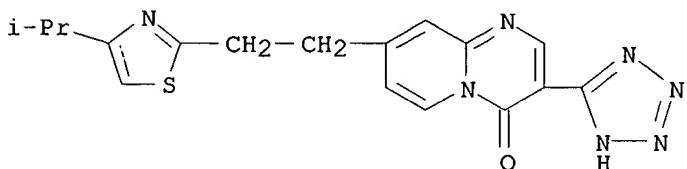
RN 337904-52-6 CAPLUS

CN 4H-Pyrido[1,2-a]pyrimidin-4-one, 8-[(4-(1-methylethyl)-2-thiazolyl)methoxy]-3-(1H-tetrazol-5-yl)- (9CI) (CA INDEX NAME)



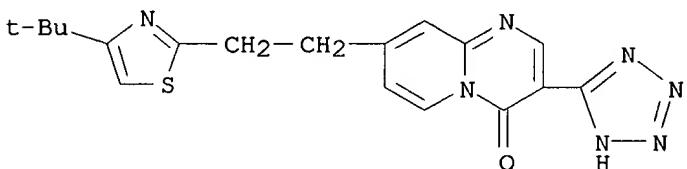
RN 337904-53-7 CAPLUS

CN 4H-Pyrido[1,2-a]pyrimidin-4-one, 8-[2-[4-(1-methylethyl)-2-thiazolyl]ethyl]-3-(1H-tetrazol-5-yl)- (9CI) (CA INDEX NAME)



RN 337904-56-0 CAPLUS

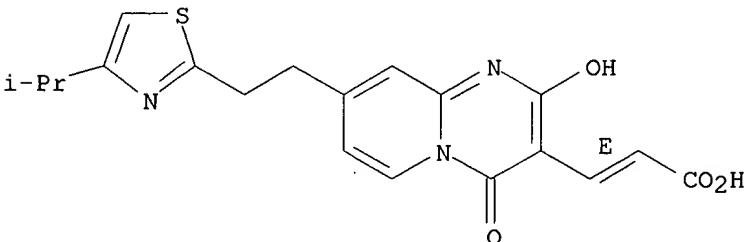
CN 4H-Pyrido[1,2-a]pyrimidin-4-one, 8-[2-[4-(1,1-dimethylethyl)-2-thiazolyl]ethyl]-3-(1H-tetrazol-5-yl)- (9CI) (CA INDEX NAME)



RN 337904-57-1 CAPLUS

CN 2-Propenoic acid, 3-[2-hydroxy-8-[2-[4-(1-methylethyl)-2-thiazolyl]ethyl]-4-oxo-4H-pyrido[1,2-a]pyrimidin-3-yl]-, (2E)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

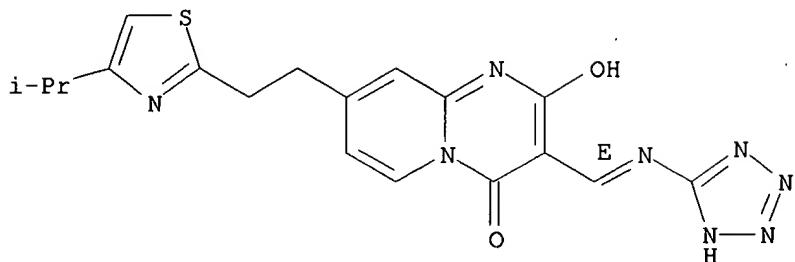


RN 337904-58-2 CAPLUS

CN 4H-Pyrido[1,2-a]pyrimidin-4-one, 2-hydroxy-8-[2-[4-(1-methylethyl)-2-thiazolyl]ethyl]-3-[(E)-(1H-tetrazol-5-ylimino)methyl]- (9CI) (CA INDEX)

NAME)

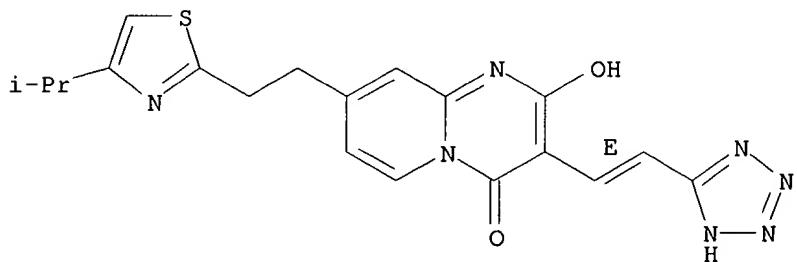
Double bond geometry as shown.



RN 337904-59-3 CAPLUS

CN 4H-Pyrido[1,2-a]pyrimidin-4-one, 2-hydroxy-8-[2-[4-(1-methylethyl)-2-thiazolyl]ethyl]-3-[(1E)-2-(1H-tetrazol-5-yl)ethenyl]- (9CI) (CA INDEX NAME)

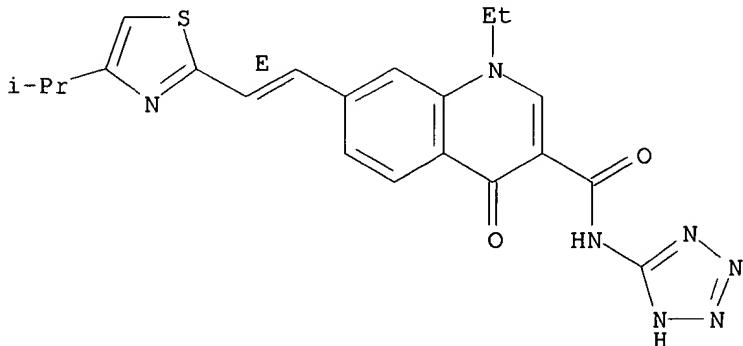
Double bond geometry as shown.



RN 337904-60-6 CAPLUS

CN 3-Quinolinecarboxamide, 1-ethyl-1,4-dihydro-7-[(1E)-2-[4-(1-methylethyl)-2-thiazolyl]ethenyl]-4-oxo-N-1H-tetrazol-5-yl- (9CI) (CA INDEX NAME)

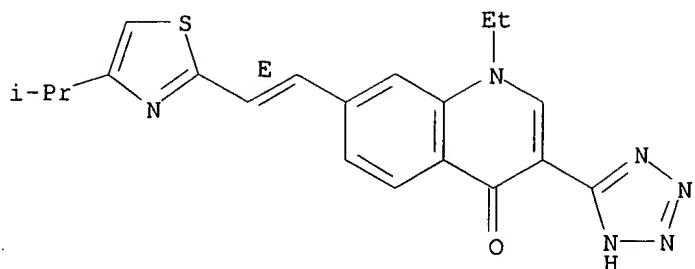
Double bond geometry as shown.



RN 337904-61-7 CAPLUS

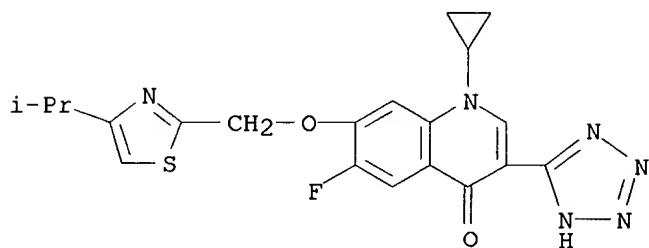
CN 4(1H)-Quinolinone, 1-ethyl-7-[(1E)-2-[4-(1-methylethyl)-2-thiazolyl]ethenyl]-3-(1H-tetrazol-5-yl)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



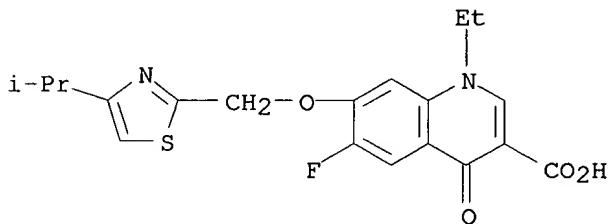
RN 337904-62-8 CAPLUS

CN 4(1H)-Quinolinone, 1-cyclopropyl-6-fluoro-7-[[4-(1-methylethyl)-2-thiazolyl]methoxy]-3-(1H-tetrazol-5-yl)- (9CI) (CA INDEX NAME)



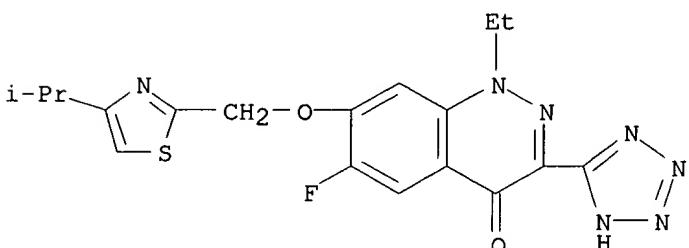
RN 337904-63-9 CAPLUS

CN 3-Quinolincarboxylic acid, 1-ethyl-6-fluoro-1,4-dihydro-7-[[4-(1-methylethyl)-2-thiazolyl]methoxy]-4-oxo- (9CI) (CA INDEX NAME)



RN 337904-64-0 CAPLUS

CN 4(1H)-Cinnolinone, 1-ethyl-6-fluoro-7-[[4-(1-methylethyl)-2-thiazolyl]methoxy]-3-(1H-tetrazol-5-yl)- (9CI) (CA INDEX NAME)



09/842,234

RE.CNT 57 THERE ARE 57 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

09/842,234

=> d his

(FILE 'HOME' ENTERED AT 09:41:47 ON 11 FEB 2002)

FILE 'REGISTRY' ENTERED AT 09:42:01 ON 11 FEB 2002

L1 STRUCTURE UPLOADED
L2 1 S L1 SSS SAM
L3 22 S L1 SSS FUL

FILE 'CAPLUS' ENTERED AT 09:43:49 ON 11 FEB 2002

L4 2 S L3

FILE 'CAOLD' ENTERED AT 09:44:30 ON 11 FEB 2002

=> s 13

L5 0 L3

=> log y

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.32	150.44

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-1.24

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